

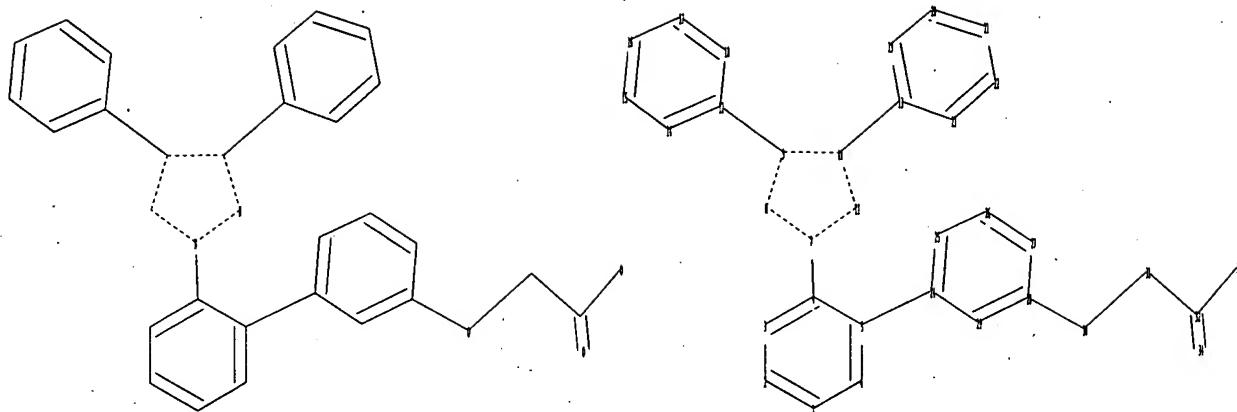
EAST Search History

Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
L1	796	(548/252):CCLS.	US-PGP UB; USPAT; USOCR ; EPO; JPO; DERWE NT; IBM_T DB	OR	OFF	2006/12/10 19:13
L2	577	(548/341.1).CCLS.	US-PGP UB; USPAT; USOCR ; EPO; JPO; DERWE NT; IBM_T DB	OR	OFF	2006/12/10 19:13
L3	828	(548/235).CCLS.	US-PGP UB; USPAT; USOCR ; EPO; JPO; DERWE NT; IBM_T DB	OR	OFF	2006/12/10 19:14

EAST Search History

L4	21	L1 AND L2	US-PGP UB; USPAT; USOCR ; EPO; JPO; DERWE NT; IBM_T DB	OR	ON	2006/12/10 19:14
L5	5	L3 AND L4	US-PGP UB; USPAT; USOCR ; EPO; JPO; DERWE NT; IBM_T DB	OR	ON	2006/12/10 19:15

10726076



chain nodes :

30 31 32 33 34

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23
24 25 26 27 28 29

chain bonds :

4-7 5-24 9-12 10-13 28-30 30-31 31-32 32-33 32-34

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 7-8 7-11 8-9 9-10 10-11 12-14 12-18 13-19
13-23 14-15 15-16 16-17 17-18 19-20 20-21 21-22 22-23 24-25 24-29 25-26
26-27 27-28 28-29

exact/norm bonds :

4-7 7-8 7-11 8-9 9-10 10-11 28-30 30-31 32-33 32-34

exact bonds :

5-24 9-12 10-13 31-32

normalized bonds :

1-2 1-6 2-3 3-4 4-5 5-6 12-14 12-18 13-19 13-23 14-15 15-16 16-17
17-18 19-20 20-21 21-22 22-23 24-25 24-29 25-26 26-27 27-28 28-29

isolated ring systems :

containing 1 : 12 : 13 :

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom
11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom
20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:Atom 26:Atom 27:Atom 28:Atom
29:Atom 30:CLASS 31:CLASS 32:CLASS 33:CLASS 34:CLASS

L1 STRUCTURE UPLOADED

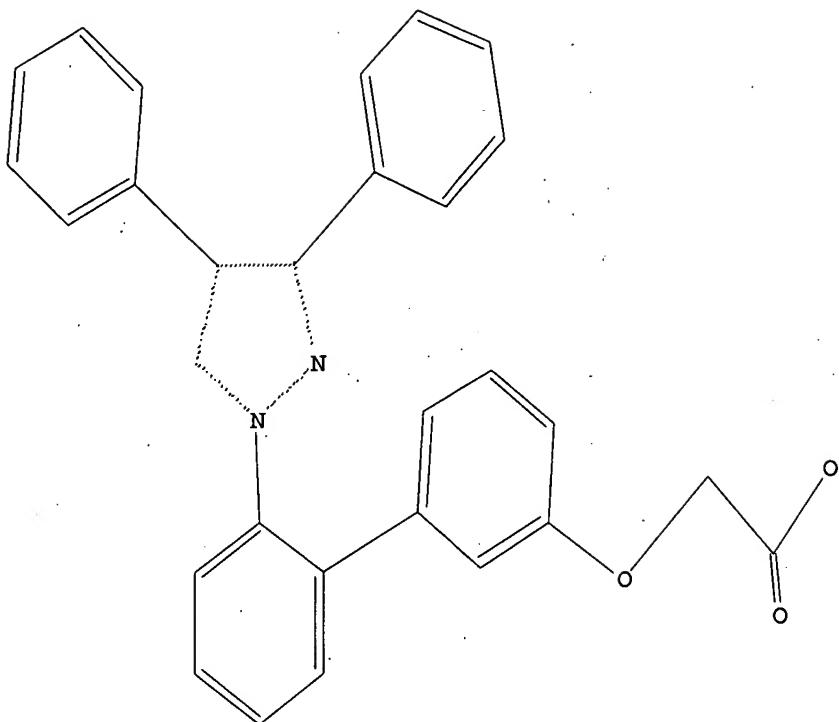
=> D

L1 HAS NO ANSWERS

L1 STR

SAEED

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Structure attributes must be viewed using STN Express query preparation.

=> S L4

L4 NOT FOUND

The L-number entered has not been defined in this session, or it has been deleted. To see the L-numbers currently defined in this session, enter DISPLAY HISTORY at an arrow prompt (=>).

=> S L1

SAMPLE SEARCH INITIATED 19:00:31 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 0 TO ITERATE

100.0% PROCESSED 0 ITERATIONS

SEARCH TIME: 00.00.01

0 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 0 TO 0

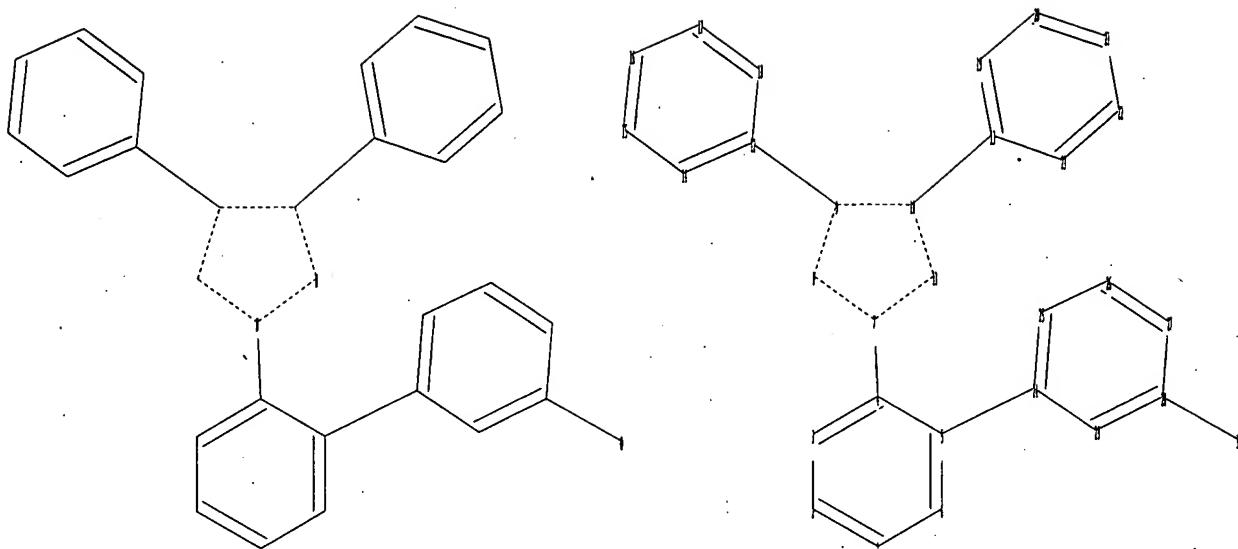
PROJECTED ANSWERS: 0 TO 0

L2 0 SEA SSS SAM L1

=>
Uploading C:\Program Files\Stnexp\Queries\107260761.str

SAEED

10726076.



chain nodes :

30

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23
24 25 26 27 28 29

chain bonds :

4-7 5-24 9-12 10-13 28-30

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 7-8 7-11 8-9 9-10 10-11 12-14 12-18 13-19
13-23 14-15 15-16 16-17 17-18 19-20 20-21 21-22 22-23 24-25 24-29 25-26
26-27 27-28 28-29

exact/norm bonds :

4-7 7-8 7-11 8-9 9-10 10-11 28-30

exact bonds :

5-24 9-12 10-13

normalized bonds :

1-2 1-6 2-3 3-4 4-5 5-6 12-14 12-18 13-19 13-23 14-15 15-16 16-17
17-18 19-20 20-21 21-22 22-23 24-25 24-29 25-26 26-27 27-28 28-29

isolated ring systems :

containing 1 : 12 : 13 :

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom
11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom
20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:Atom 26:Atom 27:Atom 28:Atom
29:Atom 30:CLASS

L3 STRUCTURE UPLOADED

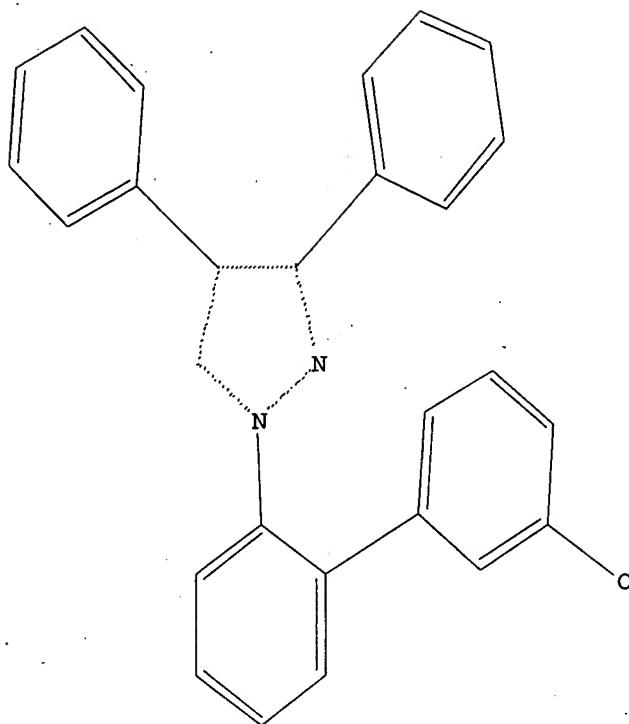
=> D

L3 HAS NO ANSWERS

L3 STR

SAEED

10726076



Structure attributes must be viewed using STN Express query preparation.

=> S L3

SAMPLE SEARCH INITIATED 19:02:06 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 1 TO ITERATE

100.0% PROCESSED 1 ITERATIONS 0 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 1 TO 80

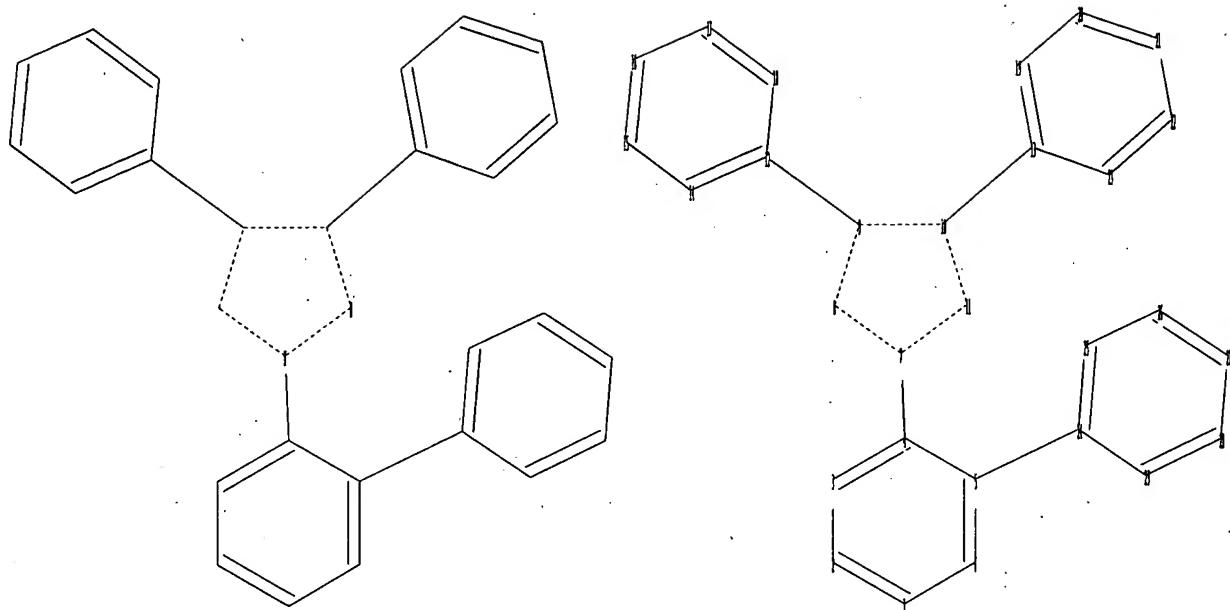
PROJECTED ANSWERS: 0 TO 0

L4 0 SEA SSS SAM L3

=>
Uploading C:\Program Files\Stnexp\Queries\107260762.str

SAEED

10726076



ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23
24 25 26 27 28 29

chain bonds :

4-7 5-24 9-12 10-13

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 7-8 7-11 8-9 9-10 10-11 12-14 12-18 13-19
13-23 14-15 15-16 16-17 17-18 19-20 20-21 21-22 22-23 24-25 24-29 25-26
26-27 27-28 28-29

exact/norm bonds :

4-7 7-8 7-11 8-9 9-10 10-11

exact bonds :

5-24 9-12 10-13

normalized bonds :

1-2 1-6 2-3 3-4 4-5 5-6 12-14 12-18 13-19 13-23 14-15 15-16 16-17
17-18 19-20 20-21 21-22 22-23 24-25 24-29 25-26 26-27 27-28 28-29

isolated ring systems :

containing 1 : 12 : 13 :

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom
11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom
20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:Atom 26:Atom 27:Atom 28:Atom
29:Atom

L5 STRUCTURE UPLOADED

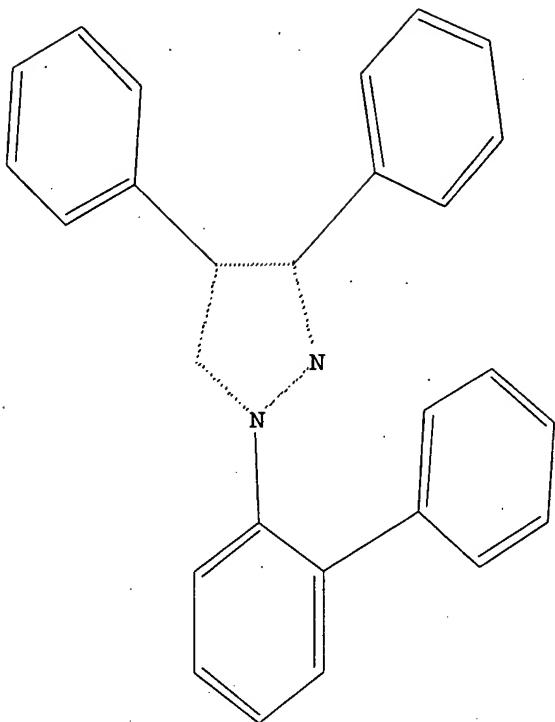
=> D

L5 HAS NO ANSWERS

L5 STR

SAEED

10726076



Structure attributes must be viewed using STN Express query preparation.

=> S L5
SAMPLE SEARCH INITIATED 19:03:28 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 7 TO ITERATE

100.0% PROCESSED 7 ITERATIONS 2 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 7 TO 298
PROJECTED ANSWERS: 2 TO 124

L6 2 SEA SSS SAM L5

=> S L5 FULL
FULL SEARCH INITIATED 19:03:36 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 109 TO ITERATE

100.0% PROCESSED 109 ITERATIONS 24 ANSWERS
SEARCH TIME: 00.00.01

L7 24 SEA SSS FUL L5

=> FILE CAPLUS
COST IN U.S. DOLLARS
FULL ESTIMATED COST SINCE FILE TOTAL
ENTRY SESSION
169.14 169.35

SAEED

10726076

FILE 'CAPLUS' ENTERED AT 19:03:41 ON 10 DEC 2006
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FILE LAST UPDATED: 8 Dec 2006 (20061208/ED)

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=> S L6
L8 1 L6

=> D IBIB ABS HITSTR TOT

SAEED

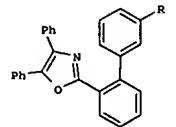
10726076

L8 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2006 ACS on STN
 ACCESSION NUMBER: 2000-725467 CAPLUS
 DOCUMENT NUMBER: 133:296436
 TITLE: Heterocyclibiphenyl α P2 inhibitors
 INVENTOR(S): Robl, Jeffrey A.; Sulsky, Richard B.; Magnin, David R.
 PATENT ASSIGNEE(S): Bristol-Myers Squibb Co., USA
 SOURCE: PCT Int. Appl., 206 pp.
 CODEN: PIKXD2
 DOCUMENT TYPE: Patent
 LANGUAGE: English
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2000059506	A1	20001012	WO 2000-US7417	20000320
W: AB, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CR, CU, DE, DK, DM, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MN, MM, MX, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, VN, YU, ZA, ZM				
RM: GH, GM, KE, LS, MM, SD, SL, SZ, TZ, UG, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BP, BJ, CF, CG, CI, CM, GN, GW, MR, NE, SN, TD, TO				
US 6548529	B1	20030415	US 2000-519079	20000320
CA 2366871	AA	20001012	CA 2000-2366871	20000320
BR 200009563	AA	20020115	BR 2000-9563	20000320
EP 1181014	A1	20020227	EP 2000-91817	20000320
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO				
TR 200102874	T3	20020321	TR 2001-2874	20000320
HU 200202251	A2	20021120	HU 2002-2251	20000320
JP 2002541106	I2	20021203	JP 2000-609070	20000320
ES 20010504	A	20021216	EE 2001-504	20000320
NZ 513493	A	20040226	NZ 2000-513493	20000320
ZA 200106856	A	20021120	ZA 2001-6956	20010820
LT 4921	B	20020625	LT 2001-92	20010925
BG 105968	B	20020531	BG 2001-105968	20011002
NO 200104623	A	20011004	NO 2001-4823	20011004
LV 12782	B	20020620	LV 2001-155	20011102
US 2003199563	A1	20031023	US 2002-321137	20021217
US 6927227	B2	20050809		
PRIORITY APPLN. INFO.:			US 1999-127745P	P 19990405
			US 2000-519079	A3 20000306
			WO 2000-US7417	W 20000320

OTHER SOURCE(S): MARPAT 133:296436
 GI

L8 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



AB α P2 inhibiting biphenyls substituted in the 2-position by a substituted 5-membered heterocycle and in the 3'-position by a carboxyalkyl, carboxymethoxy, carboxymethylamino, or 5-tetrazoylmethyl group, were prepared. The compds. are useful for treating diabetes and related diseases, especially Type II diabetes (no data) and may be used in combination with another antidiabetic agent such as metformin, gliburide, troglitazone and/or insulin. Thus, 2-BuC₆H₄CO₂H was treated with benzoin and the resulting keto ester cyclized to give 2-(2-bromophenyl)-4,5-diphenyloxazole which was coupled with 3-OCH₂C₆H₄B(OH)₂ to give the biphenyl derivative I [R = CHO]. Reduction of the formyl group, chlorination, and reaction with NaCN gave I [R = CH₂CN] which was cyclized with Me₃SnN₃ to give I [R = 5-tetrazoylmethyl].

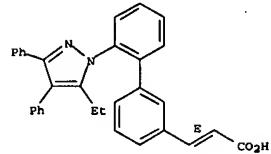
IT 300657-09-4P 300657-14-1P

RL: SPR (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses); (preparation of heterocyclibiphenyl derivs. as α P2 inhibitors)

RN 300657-09-4 CAPLUS

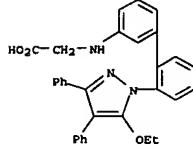
CN 2-Propenoic acid, 3-[2'-(5-ethyl-3,4-diphenyl-1H-pyrazol-1-yl)[1,1'-biphenyl]-3-yl]-, (2E)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 300657-14-1 CAPLUS
 CN Glycine, N-[2'-(5-ethoxy-3,4-diphenyl-1H-pyrazol-1-yl)[1,1'-biphenyl]-3-yl]- (9CI) (CA INDEX NAME)

L8 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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10726076

=> LOGOFF

ALL L# QUERIES AND ANSWER SETS ARE DELETED AT LOGOFF

LOGOFF? (Y)/N/HOLD:Y

COST IN U.S. DOLLARS

FULL ESTIMATED COST

SINCE FILE ENTRY	TOTAL SESSION
5.57	174.92

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE ENTRY	TOTAL SESSION
-0.75	-0.75

CA SUBSCRIBER PRICE

STN INTERNATIONAL LOGOFF AT 19:04:30 ON 10 DEC 2006

SAEED